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CENTRAL FAX CENTER
MAY 2 7 2010

Claim Amendments.

1. (currently amended): A compound of formula I:

or a pharmaceutically acceptable derivative salt, ester, salt of an ester, stereoisomer, enantiomer, isotope, or tautomer thereof, wherein:

ring A is optionally substituted and is

each R1 and R2 is independently H, alkyl, or fluoroalkyl;

R³ is H, alkyl, fluoroalkyl, aralkyl, carbocyclylalkyl, heterocyclyl, carbocyclyl, heterocyclylalkyl, aryl, heteroaryl, heteroaralkyl, -C(O)R, -OR,

 $-(CH_2)_{1-6}OR$, $-(CH_2)_{1-6}N(R)_2$, $-N(R)_2$, or -C(H)(OR)R;

 R^4 is H, alkyl, fluoroalkyl, -CO₂R, -CON(R)₂, carbocyclyl, carbocyclylalkyl, heteroaryl, or heterocyclyl;

 R^5 is $-OR^7$ or $-NR^8R^9$;

 R^6 is -C(O)R, -C(S)R, -C=C-C(O)R, -SR, $-S-W-OR^7$, M, or Y;

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R<sup>7</sup> is R°, -C(O)R, -C(O)N(R)<sub>2</sub>, -C(O)OR, -(CH<sub>2</sub>)<sub>1-6</sub>-C(O)R, -PO<sub>3</sub>M<sub>x</sub>,
-P(O)(alkyl)OM', -(PO<sub>3</sub>)<sub>2</sub>M<sub>y</sub>, carbocyclyl, aryl, heterocyclyl, heteroaryl,
carbocyclylalkyl, aralkyl, heterocyclylalkyl, or heteroaralkyl, or a tumor-targeting
moiety;
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x is 1 or 2;
y is 1, 2 or 3;
each M is independently H, Li, Na, K, Mg, Ca, Mn, Co, Ni, Zn, or alkyl;
M' is H, Li, Na, K, or alkyl;
R⁸ is H or alkyl;

 R^9 is H, alkyl, -C(O)R, $-C(O)N(R)_2$, -C(O)OR, $-SO_2R$, $-SO_2N(R)_2$, carbocyclyl, aryl, heterocyclyl, heterocyclylalkyl, aralkyl, heterocyclylalkyl, or heteroaralkyl or a tumor targeting moiety;

each R^a and R^b is independently H, OR°, alkyl, or fluoroalkyl; each R° and R^d is independently H, alkyl, or fluoroalkyl; n is 0-4;

W is alkylene, arylene, heteroarylene, carbocyclylene, or heterocyclylene; R° is H or alkyl; and

R is R°, carbocyclyl, aryl, heterocyclyl, heteroaryl, carbocyclylalkyl, aralkyl, heterocyclylalkyl, or heteroaralkyl.

- 2. (previously presented) The compound of claim 1, wherein R⁶ is Y or -SR.
- 3. (cancelled).

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- 4. (currently amended) The compound of claim 1, wherein:
 - i) R¹, R² and R⁴ are independently H, C₁₋₆ alkyl or fluoro(C₁₋₆ alkyl);
 - ii) R^3 is H, alkyl, fluoroalkyl, -(CH₂)₁₋₆OR, -(CH₂)₁₋₆N(R)₂,
- -NR°C(O)R, -C(O)R, -C(H)(OR)R, aralkyl, heterocyclyl, heterocyclylalkyl, heteroaryl, or heteroaralkyl;
 - iii) R^6 is -C=C-C(O)R, -SR, -S-W-OR⁷, M-or Y:
- iv) R⁷ is H, alkyl, -C(O)R, -PO₃M_x, -(PO₃)₂M_y, -P(O)(alkyl)OM',
 -C(O)N(R)₂, or -C(O)OR, or a turnor targeting moiety; or R⁹ is H, alkyl, -C(O)R,
 -C(O)N(R)₂, -C(O)OR, -SO₂R, 5-membered heterocyclyl, or a 5-membered
 - v) n is 1.

heteroaralkyl, or a tumor targeting moiety; and

- 5. (previously presented) The compound of claim 4, wherein R is R°, carbocyclyl, aryl, heterocyclyl, aralkyl, heterocyclylalkyl or heteroaralkyl.
- 6. (previously presented) The compound of claim 5, wherein R^0 is H or C_{1-6} alkyl optionally substituted with halo, hydroxy or amino.
- 7. (previously presented) The compound of claim 4, wherein:
- i) ring A is optionally substituted with $-NH_2$, alkyl, $-OC(O)R^{\dagger}$, halo, $-OR^{\dagger}$, $-CF_3$, $-OCF_3$, $-SCF_3$, $-SR^{\dagger}$, $-R^{\dagger}$, $-NR^{\dagger}C(O)R^{\dagger}$, $-CO_2R^{\dagger}$, $-NO_2$, $-N(R^{\dagger})_2$, -CN, $-C(O)R^{\dagger}$, $-C(O)N(R^{\dagger})_2$, $-SO_2N(R^{\dagger})_2$, $-NR^{\dagger}CO_2R^{\dagger}$, $-C(O)C(O)R^{\dagger}$, $-OC(O)N(R^{\dagger})_2$, $-S(O)_kR^{\dagger}$, $-C(O)CH_2C(O)R^{\dagger}$, $-NR^{\dagger}SO_2R^{\dagger}$, or $-C(=S)N(R^{\dagger})_2$; and R^{\dagger} is 3-6 membered unsubstituted cycloalkyl, phenyl, benzyl, naphthyl, pyridyl, or C_{1-6} alkyl optionally substituted with halo;
 - ii) R^3 is H, C_{1-6} alkyl, -(CH₂)₁₋₆OR° or -CH(OR°)R°;

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- iii) R^6 is -C=C-C(O)R, -SR, -S-W-OR⁷ or Y; and
- iv) R⁸ is H or C₁₋₆ unsubstituted alkyl.
- 8. (currently amended) The compound of claim 7, wherein R^7 or R^9 is H, a polysaccharide, $\frac{\{C(O)CH(R)N(R)\}_{2,3}-R$, an antibody, or

- 9. (currently amended) The compound of claim 7, wherein:
 - i) R¹, R² and R⁴ are independently H, methyl, ethyl, -CH₂F, -CHF₂, or -CF₃;
 - ii) R³ is H, methyl, ethyl, -CH(OH)CH₃, -CH₂OH, or -CH₂CH₂OH;
 - iii) R^6 is -S-(heterocyclylalkyl), (-S-(unsubstituted C_{1-6} alkyl), Y,

- iv) R⁸ is H, methyl, or ethyl; and
- v) R⁷ is H, methyl, ethyl, -C(O)Me, -C(O)Et, -C(O)NMe₂, -C(O)-p-OMephenyl, -C(O)O-phenyl, -PO₃H₂, -P(O)(OMe)₂, -P(O)(OMe)OH, -P(O)(Me)OH, -P(O)(OH)OP(O)(OH)(OH), or R¹¹; and R¹¹ is selected from the group consisting of:

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, and an antibody; or

R⁹ is H, methyl, ethyl, R¹¹,

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10. (currently amended) The compound of claim1, wherein said compound is selected from the group consisting of the compounds of:

(1) formula Ha:

where R^3 and R^4 are independently H or alkyl, R^6 is -SR, R^7 is R° , and R^* can be the same or different and is selected from the group consisting of alkyl and NH_{24}

(2) formulae III 13 to III 18,:

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and

(3) formulae IV 13 to IV 18:

11. (previously presented) A pharmaceutical composition comprising a compound of claim 1 and a pharmaceutically acceptable carrier.

12.-22. (cancelled).

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23. (currently amended) A compound of the formula:

$$\begin{array}{c} H_2N \\ \end{array}$$

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(d)·

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(f)

(g)

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(h)

(i)

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(j)

(k)

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(1)

(m)

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(n)

(o)

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or a pharmaceutically acceptable derivative salt, ester, salt of an ester, stereoisomer, enantiomer, isotope, or tautomer thereof.

24. (previously presented) The compound of claim 23, wherein the compound is:

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COLMAN SUDOL SAPONE

25. (currently amended) A compound of the formula

$$\begin{pmatrix}
R^{2} \\
\Pi
\end{pmatrix}$$

$$\begin{pmatrix}
R^{2} \\
\Pi
\end{pmatrix}$$

$$\begin{pmatrix}
R^{3} \\
R^{4}
\end{pmatrix}$$

$$\begin{pmatrix}
R^{3} \\
R^{6}
\end{pmatrix}$$

$$\begin{pmatrix}
R^{3} \\
R^{6}
\end{pmatrix}$$

or a pharmaceutically acceptable derivative salt, ester, salt of an ester, stereoisomer, enantiomer, isotope, or tautomer thereof, wherein:

- (a) R³ and R⁴ may each be the same or different to the extent they occur more than once in the compound and are independently H or alkyl;
- (b) R⁷ may be the same or different to the extent it occurs more than once in the compound and is independently R° or -C(O)R, where R° is H or alkyl and R is R°, carbocyclyl, aryl, heterocyclyl, heterocyclyl, carbocyclylalkyl, aralkyl, heterocyclylalkyl, or heteroaralkyl;
- (c) R^x may be the same or different to the extent it occurs more than once in the compound and is independently alkyl or NH₂;

(d)
$$R^6$$
 is $-SR$, $-C(O)R$,

or

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$$\begin{pmatrix}
R^{x} \end{pmatrix}_{11}^{N} \\
R^{4} \\
R^{3} \\
R^{3}$$

; and

(e) n is 0, 1, 2, or 3.